Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA N\Data\BN072518\
Data File : BN002139.D
Acq On : 25 Jul 2018 17:07
Operator : JU/SJ
Sample : SSTDICV040
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument : BNA_N
ClientSampleId : ICVBN072518

Quant Time: Jul 25 20:09:10 2018
Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\8270-BN072518.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Jul 25 20:08:02 2018
Response via : Initial Calibration

Min. RRF : 0.000 Min. Rel. Area : 50% Max. R.T. Dev 0.50min
Max. RRF Dev : 25% Max. Rel. Area : 150%

		Compound	Amount	Calc.	%Dev	Area%	Dev(min)
1	I	1,4-Dichlorobenzene-d4	20.000	20.000	0.0	93	0.00
2		1,4-Dioxane	40.000	39.286	1.8	91	0.00
3		Pyridine	40.000	39.927	0.2	90	0.00
4		n-Nitrosodimethylamine	40.000	36.999	7.5	89	0.00
5	S	2-Fluorophenol	80.000	77.827	2.7	90	0.00
6		Aniline	40.000	37.926	5.2	88	0.00
7	S	Phenol-d6	80.000	76.173	4.8	87	0.00
8		2-Chlorophenol	40.000	38.251	4.4	88	0.00
9		Benzaldehyde	40.000	38.500	3.8	90	0.00
10	C	Phenol	40.000	38.094	4.8	88	0.00
11		bis(2-Chloroethyl)ether	40.000	38.504	3.7	90	0.00
12		1,3-Dichlorobenzene	40.000	38.388	4.0	89	0.00
13	C	1,4-Dichlorobenzene	40.000	38.231	4.4	89	0.00
14		1,2-Dichlorobenzene	40.000	37.736	5.7	88	0.00
15		Benzyl Alcohol	40.000	36.978	7.6	86	0.00
16		2,2'-oxybis(1-Chloropropane	40.000	37.052	7.4	88	0.00
17		2-Methylphenol	40.000	37.530	6.2	87	0.00
18		Hexachloroethane	40.000	38.529	3.7	89	0.00
19	P	n-Nitroso-di-n-propylamine	40.000	36.078	9.8	87	0.00
20		3+4-Methylphenols	40.000	37.430	6.4	88	0.00
21	I	Naphthalene-d8	20.000	20.000	0.0	92	0.00
22		Acetophenone	40.000	38.149	4.6	88	0.00
23	S	Nitrobenzene-d5	80.000	77.401	3.2	88	0.00
24		Nitrobenzene	40.000	38.496	3.8	88	0.00
25		Isophorone	40.000	37.099	7.3	87	0.00
26	C	2-Nitrophenol	40.000	38.700	3.2	88	0.00
27		2,4-Dimethylphenol	40.000	38.142	4.6	87	0.00
28		bis(2-Chloroethoxy)methane	40.000	37.673	5.8	88	0.00
29	C	2,4-Dichlorophenol	40.000	38.255	4.4	88	0.00
30		1,2,4-Trichlorobenzene	40.000	38.309	4.2	88	0.00
31		Naphthalene	40.000	38.040	4.9	88	0.00
32		Benzoic acid	40.000	36.000	10.0	86	0.00
33		4-Chloroaniline	40.000	37.505	6.2	87	0.00
34	C	Hexachlorobutadiene	40.000	38.640	3.4	89	0.00
35		Caprolactam	40.000	35.208	12.0	85	-0.01
36	C	4-Chloro-3-methylphenol	40.000	36.499	8.8	86	0.00
37		2-Methylnaphthalene	40.000	37.335	6.7	88	0.00
38	I	Acenaphthene-d10	20.000	20.000	0.0	91	0.00
39		1,2,4,5-Tetrachlorobenzene	40.000	39.987	0.0	87	0.00
40	P	Hexachlorocyclopentadiene	40.000	42.111	-5.3	87	0.00
41	S	2,4,6-Tribromophenol	80.000	73.975	7.5	85	0.00
42	C	2,4,6-Trichlorophenol	40.000	39.958	0.1	88	0.00
43		2,4,5-Trichlorophenol	40.000	39.287	1.8	87	0.00
44	S	2-Fluorobiphenyl	80.000	79.889	0.1	88	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA N\Data\BN072518\
Data File : BN002139.D
Acq On : 25 Jul 2018 17:07
Operator : JU/SJ
Sample : SSTDICV040
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument : BNA_N
ClientSampleId : ICVBN072518

Quant Time: Jul 25 20:09:10 2018
Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\8270-BN072518.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Jul 25 20:08:02 2018
Response via : Initial Calibration

Min. RRF : 0.000 Min. Rel. Area : 50% Max. R.T. Dev 0.50min
Max. RRF Dev : 25% Max. Rel. Area : 150%

		Compound	Amount	Calc.	%Dev	Area%	Dev(min)
45		1,1'-Biphenyl	40.000	39.369	1.6	88	0.00
46		2-Chloronaphthalene	40.000	39.734	0.7	88	0.00
47		2-Nitroaniline	40.000	38.437	3.9	86	0.00
48		Acenaphthylene	40.000	38.708	3.2	87	0.00
49		Dimethylphthalate	40.000	37.447	6.4	87	0.00
50		2,6-Dinitrotoluene	40.000	38.631	3.4	88	0.00
51	C	Acenaphthene	40.000	38.939	2.7	88	0.00
52		3-Nitroaniline	40.000	37.762	5.6	85	0.00
53	P	2,4-Dinitrophenol	40.000	36.144	9.6	84	0.00
54		Dibenzofuran	40.000	37.952	5.1	86	0.00
55	P	4-Nitrophenol	40.000	37.760	5.6	85	0.00
56		2,4-Dinitrotoluene	40.000	37.428	6.4	85	0.00
57		Fluorene	40.000	37.679	5.8	86	0.00
58		2,3,4,6-Tetrachlorophenol	40.000	37.021	7.4	84	0.00
59		Diethylphthalate	40.000	36.502	8.7	85	0.00
60		4-Chlorophenyl-phenylether	40.000	37.456	6.4	86	0.00
61		4-Nitroaniline	40.000	37.659	5.9	86	0.00
62		Azobenzene	40.000	37.526	6.2	86	0.00
63	I	Phenanthrene-d10	20.000	20.000	0.0	90	0.00
64		4,6-Dinitro-2-methylphenol	40.000	39.359	1.6	84	0.00
65	c	n-Nitrosodiphenylamine	40.000	38.466	3.8	85	0.00
66		4-Bromophenyl-phenylether	40.000	38.832	2.9	86	0.00
67		Hexachlorobenzene	40.000	38.070	4.8	85	0.00
68		Atrazine	40.000	38.516	3.7	86	0.00
69	C	Pentachlorophenol	40.000	38.642	3.4	83	0.00
70		Phenanthrene	40.000	37.662	5.8	85	0.00
71		Anthracene	40.000	38.063	4.8	85	0.00
72		Carbazole	40.000	37.597	6.0	84	0.00
73		Di-n-butylphthalate	40.000	37.918	5.2	84	0.00
74	C	Fluoranthene	40.000	37.736	5.7	84	0.00
75	I	Chrysene-d12	20.000	20.000	0.0	90	0.00
76		Benzidine	40.000	36.271	9.3	91	0.00
77		Pyrene	40.000	36.604	8.5	85	0.00
78	S	Terphenyl-d14	80.000	73.341	8.3	86	0.00
79		Butylbenzylphthalate	40.000	37.251	6.9	84	0.01
80		Benzo(a)anthracene	40.000	37.103	7.2	84	0.01
81		3,3'-Dichlorobenzidine	40.000	37.832	5.4	83	0.01
82		Chrysene	40.000	37.341	6.6	84	0.01
83		Bis(2-ethylhexyl)phthalate	40.000	38.337	4.2	86	0.01
84	c	Di-n-octyl phthalate	40.000	39.430	1.4	85	0.02
85		Indeno(1,2,3-cd)pyrene	40.000	41.629	-4.1	83	0.02
86	I	Perylene-d12	20.000	20.000	0.0	89	0.02
87		Benzo(b)fluoranthene	40.000	37.813	5.5	84	0.02

Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA N\Data\BN072518\
Data File : BN002139.D
Acq On : 25 Jul 2018 17:07
Operator : JU/SJ
Sample : SSTDICV040
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument : BNA_N
ClientSampleId : ICVBN072518

Quant Time: Jul 25 20:09:10 2018
Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\8270-BN072518.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Jul 25 20:08:02 2018
Response via : Initial Calibration

Min. RRF : 0.000 Min. Rel. Area : 50% Max. R.T. Dev 0.50min
Max. RRF Dev : 25% Max. Rel. Area : 150%

	Compound	Amount	Calc.	%Dev	Area%	Dev(min)
88	Benzo(k)fluoranthene	40.000	36.998	7.5	84	0.02
89 C	Benzo(a)pyrene	40.000	37.945	5.1	84	0.02
90	Dibenzo(a,h)anthracene	40.000	39.547	1.1	83	0.02
91	Benzo(g,h,i)perylene	40.000	39.879	0.3	83	0.02

(#) = Out of Range SPCC's out = 0 CCC's out = 0